# **AMENDMENTS TO THE CLAIMS**

This listing of claims will replace all prior versions, and listings, of claims in the application:

Claims 1-28 previously cancelled.

Claims 29-36 cancelled.

37. (Currently Amended) The A compound represented by formula (1) according to claim 3656, in which R5 is a phenyl group substituted at least in position 3.

38. (Currently Amended) The A compound represented by formula (11) according to claim 3656, wherein R5 is a phenyl group substituted at least in position 3 and wherein the substituent by a substituent is selected infrom the group consisting of: CHO, CN, CONH2, NO2, CF3, NH2, halogen-atom (Cl), (C1-C6) alkyl, a phenyl **group** optionally substituted; in particular by an acetyl group, by a halogen atom (Cl), by a CONH2-group or by a CN, a prop-1-ynyl group optionally substituted, in particular by a benzyloxy or tenttert-butyl carbamate, a hex-1-ynyl group optionally substituted, in particular by a CN or NH2 group, a pentyl group optionally substituted, in particular by a CONH2, a hexyl group, a piperidinyl **group** optionally substituted, in particular by a-prop-1-ynyl, benzylaminomethyl, acetamide (CH3CONH), aminomethyl, NH2CS-, 4-phenyl-1,3-thiazo-2-yl, -CONHBenzyl, -COOEthyl, -CONHiPropyl, -CONH-(CH2)n-CONH2 (wherein n representing represents a whole number from 1 to 6), -CONR'R"-group, withwherein R' and R", which are the same or different, representing represent a C1 to C6 alkyl group or a hydrogen atom, -(4-benzylpyperazin-1-yl)carbonyl, -CONH-(CH<sub>2</sub>)n-phenyl (wherein n representing represents a whole number from 1 to 6), imidazolyl, **or a** piperazinyl group optionally substituted, **in particular** by a phenyl group.

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- 39. (Currently Amended) The <u>A</u> compound represented by formula (1) according to claim <u>3656</u>, in which R5 is a phenyl group\_-substituted in positions 3 and 4, in particular by a hydrocarbon chain <u>possiblyoptionally</u> containing at least one heteroatom, <u>such as oxygen</u>, <u>like the methylenedioxy (-O-CH2-O-)said</u> chain forming a ring with the phenyl group to which it is attached.
- 40. (Currently Amended) The A compound represented by formula (11) according to claim 3656, in which R5 is the 3-pyridyl, 4-isoquinolyl, or a piperazinyl group optionally substituted, in particular in position 4. by an aryl group, such as phenyl.
- 41. (Currently Amended) The A compound represented by formula (11) according to claim 3656, in which Z represents a sulfur atom or -NR2, with in particular wherein R2 and R1 forming a form an imidazole ring of the imidazole type with R1.
- 42. (Currently Amended) The A compound represented by formula (I1) according to claim 3656, in which: Z is the oxygen atom, and/or; R<sub>7</sub> and R<sub>8</sub>, independently of each other, represent an OR<sub>10</sub> group in which R<sub>10</sub> is a-(C<sub>1</sub>-C<sub>6</sub>) alkyl group, preferably an ethyl or methyl group, advantageously methyl, and/or; R<sub>7</sub> and R<sub>8</sub> each represent an ethoxy or methoxy-group, advantageously methoxy, or one represents a hydrogen atom and the other an ethoxy or methoxy-group, advantageously methoxy, and/or; R<sub>8</sub> and R<sub>9</sub>, which are the same or different, represent the hydrogen atom, a halogen atom, a phenyl-group, a (C<sub>1</sub>-C<sub>6</sub>) alkyl-group or an OR<sub>10</sub> group in which R<sub>10</sub> is a (C<sub>1</sub>-C<sub>6</sub>) alkyl-group, preferably an ethyl or methyl group, and/or; R<sub>3</sub> and R<sub>3</sub>', which are the same or different, represent a hydrogen-atom, and/or; R<sub>1</sub> is a (C<sub>1</sub>-C<sub>6</sub>) alkyl, (C<sub>6</sub>-C<sub>18</sub>) aryl-group, such as phenyl, (C<sub>6</sub>-C<sub>18</sub>)aryl(C<sub>1</sub>-C<sub>4</sub>)alkyl-group such as benzyl, optionally substituted, or a (C<sub>1</sub>-C<sub>12</sub>)aIkyl(C<sub>6</sub>-C<sub>18</sub>)aryl group.

- 43. (Currently Amended) The  $\underline{A}$  compound represented by formula ( $\underline{I_1}$ ) according to claim  $\underline{3656}$ , in which R1 represents  $\underline{a}$ -hydrogen- $\underline{atom\ or\ a}$  (C<sub>1</sub>- $\underline{C_3C_6}$ ) alkyl, (C<sub>6</sub>-C<sub>18</sub>) aryl (G<sub>1</sub>-C<sub>4</sub>) alkyl-(for example: benzyl), or  $\underline{a}$  (C<sub>1</sub>-C<sub>12</sub>) alkyl(C<sub>6</sub>,-C<sub>18</sub>) aryl group, said  $\underline{aryl}$  group being optionally substituted.
- 44. (Currently Amended) The  $\underline{A}$  compound represented by formula ( $\underline{I_1}$ ) according to claim  $\underline{3656}$ , in which R5 is a phenyl group substituted by  $\underline{a}$  group selected from:
  - (a) one or more OR' groups, in particular methoxy or ethoxy, or
  - (b) a COR' group, in particular acetyl or aldehyde, or
  - (c) a CONR'R" group, in particular CONH2, or
  - (d) a CN group, or
  - (e) a trifluoromethyl group, or
  - (f) an alkyl group, for example methyl, or alkynyl group, for example hexynyl or propynyl, or
  - (g) an aryl or heterocycle group, in particular phenyl, furyl, pyridyl, piperidine, thiazole or thienyl, said aryl or heterocycle itself being optionally substituted by one or more groups preferably selected from groups (a)-(g).
- 45. (Currently Amended) The  $\underline{A}$  compound represented by formula ( $\underline{i}\underline{1}$ ) according to claim  $\underline{3656}$ , in which at least one  $\underline{groupof}$   $R_7$  or  $\underline{and}$   $R_8$  represents  $OR_{10}$  where  $\underline{in}$   $R_{10}$  represents a ( $C_1$ - $C_6$ ) alkyl or ( $C_3$ - $C_5$ ) cycloalkyl group.
- 46. (Currently Amended) The  $\underline{A}$  compound represented by formula ( $\underline{I_1}$ ) according to claim  $\underline{3656}$ , in which at least one group  $R_7$ , or  $R_8$  represents  $OR_{10}$  where  $\underline{in}$   $R_{10}$

represents a  $(C_1-C_6)$  alkyl or  $(C_3-C_6)$  cycloalkyl group, and at least one of **the groups**  $R_7$  and  $R_8$ , **advantageously both**, represents a methoxy group.

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47. (Currently Amended) The A compound according to claim 3656, wherein said
compound is selected from the following compoundsgroup consisting of:
3-(7,8-dimethoxy-2-oxo-2,3-dihydro-1H-1,4-benzodiazepin-5-yl)benzonitrile, 3a
7,8-dimethoxy-[5-(3-trifluoromethyl)phenyl]-1,3-dihydro-2H-1,4-benzodiazepin-2-one,
3d
3-(7,8-dimethoxy-1-methyl-2-oxo-2,3-dihydro-1H-1,4-benzodiazepin-5-yl)-benzonitrile,
4a
3-[1-(4-chlorobenzyl)-7,8-dimethoxy-2-oxo-2,3-dihydro-1H-1,4-benzodiazepin-5-yl]-
benzonitrile, 4c
3-[1-(3,4-chlorobenzyl)-7,8-dimethoxy-2-oxo-2,3-dihydro-1H-1,4-benzodiazepin-5-yl]-
benzonitrile, 4d
3-[7,8-dimethoxy-1-(4-methoxybenzyl)-2-oxo-2,3-dihydro-1H-1,4-benzodiazepin-5-yl]-
benzonitrile, 4e
3-[1-(3-chlorobenzyl)-7,8-dimethoxy-2-oxo-2,3-dihydro-1H-1,4-benzodiazepin-5-vl]-
benzonitrile, 4f
3-(7,8-dimethoxy-2-oxo-1-[3-(trfluoromethyl)benzyl]-2,3-dihydro-1H-1,4-
benzodiazepin-5-yl]-benzonitrile, 4g
3-[1-(2-chiorobenzyl)-7,8-dimethoxy-2-oxo-2,3-dihydro-1H-1,4-benzodiazepin-5-yl]-
benzonitrile, 4h
3-{7,B-dimethoxy-2-oxo-1-[4-(trifluoromethyl)benzyl]-2,3-dihydro-1H-1,4-
benzodiazepin-5-vl]-benzonitrile, 4i
3-[7,8-dimethoxy-2-oxo-1-(2-phenylethyl)-2,3-dihydro-1H-1,4-benzodiazepin-5-yl]-
benzonitrile, 4i
3-(1-ethyl-7,8-dimethoxy-2-oxo-2,3-dihydro-1H-1,4-benzodiazepin-5-yl)benzonitrile,
4k
3-(7,8-dimethoxy-l-propyl-2-oxo-2,3-dihydro-1H-1,4-benzodiazepin-5-yi)benzonitrile,
4l
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benzamide, 5j

 $5-\{3-[3-(benzyloxy)prop-l-ynyl]phenyl\}-1-ethyl-7,8-dimethoxy-1,3-dihydro-2H-1,4-benzodiazepin-2-one, \textbf{6k}$  3'-(7,8-dimethoxy-1-methyl-2-oxo-2,3-dihydro-1H-1,4-benzodiazepin-5-yl)-1,1'-biphenyl-3-carbonitrile, 61 3'-(7,8-dimethoxy-l-methyl-2-oxo-2,3-dihydro-1H-1,4-benzodiazepin-5-yl)-1,1'-biphenyl-4-carbonitrile, 6m

- 3'-(7,8-dimethoxy-1-methyl-2-oxo-2,3-dihydro-1H-1,4-benzodiazepin-5-yl)-1,1'-biphenyl-4-carboxamide, **6n**
- 3'-(7,8-dimethoxy-l-methyl-2-oxo-2,3-dihydro-1H-1,4-benzodiazepin-5-yl)-1,1'-biphenyl-3-carboxamide, **60**
- 3-[3-(3,4-dichlorobenzyl)-7,8-dimethoxy-l-methyl-2-oxo-2,3-dihydro-1H-1,4-benzodiazepin-5-yl]benzonitrile,  $7\mathbf{b}$
- 7,8-dimethoxy-1,3-dimethyl-5-(3-trifluoromethylphenyl)-1,3-dihydro-2H-1,4-benzodiazepin-2-one, 7c
- 3-(7,8-dimethoxy-1,3-dimethyl-2-oxo-2,3-dihydro-1H-1,4-benzodiazepin-5-yl)benzonitrile, **7d**
- 5-[3-(aminomethyl)phenyl]-7,8-dimethoxy-1-methyl-1,3-dihydro-2H-1,4-benzodiazepin-2-one, **8a**
- N-[3-(7,8-dimethoxy-l-methyl-2-oxo-2,3-dihydro-1 H-1,4-benzodiazepin-5-yl)benzyl]acetamide, **8b**
- 3-(7,8-dlmethoxy-1-methyl-2-oxo-2,3-dihydro-1H-1,4-benzodiazepin-5-yl)thiobenzamide, **9a**
- 7,8-dimethoxy-1-methyl-5-[3-(4-phenyl-1,3-thiazol-2-yl)phenyl]-1,3-dihydro-2H-1,4-benzodiazepin-2-one,  $\bf 9b$
- $5\hbox{-}(3\hbox{-}cyanophenyl)\hbox{-}7,8\hbox{-}dimethoxy\hbox{-}1,3\hbox{-}dihydro\hbox{-}2H\hbox{-}1,4\hbox{-}benzodiazepin\hbox{-}2\hbox{-}thione,} \textbf{10d}$
- 3-(8,9-dimethoxy-4H-imidazo[1,2-a][1,4] benzo diazepin-5-yl) benzonitrile, 11a
- 3-(8,9-dimethoxy-4H-imidazo[1,2-a)[1,4]benzodiazepin-6-yl)benzamide, 11b
- $3\hbox{-}(7,8\hbox{-}dimethoxy\hbox{-}2\hbox{-}methylamino\hbox{-}1,3\hbox{-}dihydro\hbox{-}3H\hbox{-}1,4\hbox{-}benzodiazepin\hbox{-}5\hbox{-}yl) benzonitrile,$

### 12a

- 7,8-dimethoxy-l-methyl-5-(3-pyridyl)-1,3-dihydro-1,4-benzodiazepin-2-one, 17b
- 7,8-dimethoxy-l-methyl-5-(3-nitrophenyl)-1,3-dihydro-1,4-benzodiazepin-2-one, 17c
- 5-(7,8-dimethoxy-l-methyl-2-oxo-2,3-dihydro-1H-1,4-benzodiazepin-5-yl)-2-dimethoxy-l-methyl-2-oxo-2,3-dihydro-1H-1,4-benzodiazepin-5-yl)-2-dimethoxy-l-methyl-2-oxo-2,3-dihydro-1H-1,4-benzodiazepin-5-yl)-2-dimethoxy-l-methyl-2-oxo-2,3-dihydro-1H-1,4-benzodiazepin-5-yl)-2-dimethoxy-l-methyl-2-oxo-2,3-dihydro-1H-1,4-benzodiazepin-5-yl)-2-dimethyl-2-oxo-2,3-dihydro-1H-1,4-benzodiazepin-5-yl)-2-dimethyl-2-oxo-2,3-dihydro-1H-1,4-benzodiazepin-5-yl)-2-dimethyl-2-oxo-2,3-dihydro-1H-1,4-benzodiazepin-5-yl)-2-dimethyl-2-oxo-2,3-dihydro-1H-1,4-benzodiazepin-5-yl)-2-dimethyl-2-oxo-2,3-dihydro-1H-1,4-benzodiazepin-5-yl)-2-dimethyl-2-oxo-2,3-dihydro-1H-1,4-benzodiazepin-5-yl)-2-dimethyl-2-oxo-2,3-dihydro-1H-1,4-benzodiazepin-5-yl)-2-dimethyl-2-oxo-2,3-dihydro-1H-1,4-benzodiazepin-5-yl)-2-dimethyl-2-oxo-2,3-dihydro-1H-1,4-benzodiazepin-5-yl)-2-dimethyl-2-oxo-2,3-dihydro-1H-1,4-benzodiazepin-5-yl)-2-dimethyl-2-oxo-2,3-dihydro-1H-1,4-benzodiazepin-5-yl)-2-dimethyl-2-oxo-2,3-dihydro-1H-1,4-benzodiazepin-5-yl)-2-dimethyl-2-oxo-2,3-dihydro-1H-1,4-benzodiazepin-5-yl)-2-dimethyl-2-oxo-2,3-dihydro-1H-1,4-benzodiazepin-5-yl)-2-dimethyl-2-oxo-2,3-dihydro-1H-1,4-benzodiazepin-5-yl)-2-dimethyl-2-oxo-2,3-dihydro-1H-1,4-benzodiazepin-5-yl)-2-dimethyl-2-oxo-2,3-dihydro-1H-1,4-benzodiazepin-5-yl)-2-dimethyl-2-dimethy
- benzonitrile, 17d
- 5-(3-acetylphenyl)-7,8-dimethoxy-l-methyl-1,3-dihydro-1,4-benzodiazepin-2-one, 17e
- 5-(4-isoquinolinyl)-7,8-dimethoxy-l-methyl-1,3-dihydro-1,4-benzodiazepin-2-one, 17f
- 7,8-dimethoxy-5-(3-hydroxymethylphenyl)-1-methyl-3-propyl-1,3-dihydro-2H-1,4
- benzodiazepin-2-one, 17h

5-(3-aminophenyl)-7,8-dimethoxy-1-methyl-1,3-dihydro-2H-1,4-benzodiazepin-2-one,

5-(3,4-dichlorophenyl)-7, 8-dimethoxy-1-methyl-1,3-dihydro-1,4-benzodiazepin-2-one,

7,8-dimethoxy-l-methyl-5-(3-methylphenyl)-1,3-dihydro-1,4-benzodiazepin-2-one, 17k. 5-(3-formviphenyly-7,8-dimethoxy-l-methyl-1,3-dihydro-1,4-benzodiazepin-2-one, 171 5-[3-(benzylaminomethyl)phenyl]-7,8-dimethoxy-1-methyl-1,3-dihydro-2H-1,4-

N-[3-(7,8-dimethoxy-1-methyl-2-oxo-2,3-dihydro-1H-1,4-benzodiazepin-5-yl) phenyllacetamide, 17n

7,8-dimethoxy-1-methyl-5-(3,4-methylenedioxyphenyl)-1,3-dihydro-2H-1,4benzodiazepin-2-one, 170

3-(7-hydroxy-8-methoxy-2-oxo-2,3-dihydro-111,4-benzodiazepin-5-yl)benzonitrile, 22b 3-(6-bromo-7-hydroxy-8-methoxy-2-oxo-2,3-dihydro-1H-1,4-benzodiazepin-5vl)benzonitrile, 23b

3-(9-bromo-8-hydroxy-7-methoxy-2-oxo-2,3-dihydro-1H-1,4-benzodiazepin-5yl)benzonitrile, 23d

3-(6-bromo-7,8-dimethoxy-1-methyl-2-oxo-2,3-dihydro-1H-1,4-benzodiazepin-5yl)benzonitrile, 24b

3-(7,8-dimethoxy-1-methyl-2-oxo-6-phenyl-2,3-dihydro-1H-1,4-benzodiazepin-5yl)benzonitrile, 25b

3-(7,8-dimethoxy-1-methyl-2-oxo-9-phenyl'-2,3-dihydro-1H-1,4rbenzodiazepin-5yl)benzonitrile, 25a

tert-butyl-3-[5-(cyanophenyl)-(7,8-dimethoxy-1-methyl-2-oxo-2,3-dihydro-1H-1,4benzodiazepin-9-yl)phenyl]prop-2-ynylcarbamate, 25c

methyl(2E)-3-[5-(cvanophenyl)-7,8-dimethoxy-l-methyl-2-oxo-2,3-dihydro-1H-1,4benzodiazepin-9-vl)phenyllacrylate, 25d

tert-butyl-3-[5-(cyanophenyl)-(7,8-dimethoxy-1-methyl-2-oxo-2,3-dihydro-1 H-1,4benzodiazepin-6-yl)phenyl]prop-2-ynylcarbamate, 25e

[9-(3-aminoethynyl)-7,8-dimethoxy-1-methyl-2-oxo-2,3-dihydro-H-1,4-benzodiazepin-5-yl]benzonitrile, 25f

- [6-(3-aminoethynyl)-7,8-dimethoxy-1-methyl-2-oxo-2,3-dihydro-1H-1,4-benzodiazepin-5-yl]benzonitrile, **25g**
- 3-(8-methoxy-2-oxo-2,3-dihydro-1H-1,4-benzodiazepin-5-yl)benzonitrile, 28a
- 3-(6-methoxy-2-oxo-2,3-dihydro-1H-1,4-benzodiazepin-5-yl)benzonitrile, 28b
- 3-(7-methoxy-2-oxo-2,3-dihydro-1H-1,4-benzodiazepin-5-yi)benzonitrile, 28c
- 6-methoxy-5-phenyl-1,3-dihydro-2H-1,4-benzodiazepin-2-one, 28d
- 7-methoxy-5-phenyl-1,3-dihydro-2H-1,4-benzodiazepin-2-one, 28e
- 9-bromo-7,8-dimethoxy-5-phenyl-1,3-dihydro-2H-1,4-benzodiazepin-2-one, 28f
- 3-(6,8-dimethoxy-2-oxo-2,3-dihydro-1H-1,4-benzodiazepin-5-yl)benzonitrile, 28g
- 3-(7-bromo-6,8-dimethoxy-2-oxo-2,3-dihydro-1H-1,4-benzodiazepin-5-yl)benzonitrile,

# 28h

 $3\hbox{-}(6,8\hbox{-methoxy-l-methyl-2-oxo-2,3-dihydro-1H-1,4-benzodiazepin-5-yl}) benzonitrile,$ 

## 29a

- 3-(6,8-dimethoxy-1-methyl-2-oxo-2,3-dihydro-1H-1,4-benzodiazepin-5-yl)benzonitrile, **29b**
- 3-(7-bromo-6,8-dimethoxy-1-methyl-2-oxo-2,3-dihydro-1H-1,4-benzodiazepin-5-yl)benzonitrile, **29c**
- 3-(7,8-dimethoxy-1-methyl-2-oxo-2,3-dihydro-1H-1,4-benzodiazepin-5-yl)methyl benzoate, 34a
- $3\hbox{-}(7,8\hbox{-}dimethoxy\hbox{-}1\hbox{-}methyl\hbox{-}2\hbox{-}oxo\hbox{-}2,3\hbox{-}dihydro\hbox{-}1H\hbox{-}1,4\hbox{-}benzodiazepin\hbox{-}5\hbox{-}yl)} benzoic acid,$

#### 35a

- 3-(7,8-dimethoxy-1-methyl-2-oxo-2,3-dihydro-1 H-1,4-benzodiazepin-5-yl)N-isopropylbenzamide, **36a**
- N-benzyl-3-(7,8-dimethoxy-1-methyl-2-oxo-2,3-dihydro-1H-1,4-benzodiazepin-5-yl) benzamide,  ${\bf 36b}$
- N-(6-amino-6-oxohexyl)-3-(7,8-dimethoxy-l-methyl-2-oxo-2,3-dihydro-1H-1,4-benzodiazepin-5-yl)benzamide, **36c**
- 3-(7,8-dimethoxy-l-methyl-2-oxo-2,3-dihydro-1H-1,4-benzodiazepin-5-yl)-N,N-dimethylbenzamide **36d**
- 5-(3-[(4-benzylpiperazin-1-yl)carbonyl]phenyl)7,8-dimethoxy-1-methyl-2-oxo-2,3-dihydro-1H-1,4-benzodiazepin-2-one, **36e** <u>and</u>

- 3-(7,8-dimethoxy-1-methyl-2-oxo-2,3-dihydro- 1 H-1,4-benzodiazepin-5-yl)-N-(3-phenylpropyl)benzamide, **36f**.
- 48. (Currently Amended) The The compound according to claim 36, wherein said compound is selected from the following compounds group consisting of: 3-(1-benzyl-7,8-dimethoxy-2-oxo-2,3-dihydro-1H-1,4-benzodiazepin-5-yl)benzonitrile, 4m
- 7,8-dimethoxy-1-methyl-[5-(3-trifluoromethyl)phenyl]-1,3-dihydro-2H-1,4-benzodiazepin-2-one, **4p**
- 3-(7,8-dimethoxy-1-methyl-2-oxo-2,3-dihydro-1H-1,4-benzodiazepin-5-yl)benzamide, **5a**
- 3-(6-bromo-7,8-dimethoxy-l-methyl-2-oxo-2,3-dihydro-1H-1,4-benzodiazepin-5-yl)benzamide,  ${\bf 5b}$
- tert-butyl-3-[3-(7,8-dimethoxy-1-methyl-2-oxo-2,3-dihydro-1H-1,4-benzodiazepin-5-yl)phenyl]propynylcarbamate, **6a**
- 7,8-dimethoxy-5-(3'-hex-1-ynylphenyl)-1-N-methyl-1,3-dihydro-2H-1,4-benzodiazepin-2-one,  ${\bf 6b}$
- 6-[3-(7,8-dimethoxy-1-methyl-2-oxo-2,3-dihydro-1H-1,4-benzodiazepin-5-yl)phenyl]hex-5-ynenitrile, **6d**
- 7,8-dimethoxy-5-(3'-hexylphenyl)-1-N-methyl-1,3-dihydro-2H-1,4-benzodiazepin-2-one, $\mathbf{6e}$
- $5-(4'-chloro-1,1'-biphenyl-3-yl)-7,8-dimethoxy-l-methyl-1,3-dihydro-2H-1,4-benzodiazepin-2-one, <math>\bf 6j$
- 3'-(7,8-dimethoxy-1-methyl-2-oxo-2,3-dihydro-1H-1,4-benzodiazepin-5-yl)-1,1'-biphenyl-4-carbonitrile, **6m**
- 3'-(7,8-dimethoxy-l-methyl-2-oxo-2,3-dihydro-1 H-1,4-benzodiazepin-5-yl)-1,1'-biphenyl-4-carboxamide,  $\bf 6n$
- 3-(3,4-dichlorobenzyl)-1-ethyl-7,8-dimethoxy-5-phenyl-1,3-dihydro-2H-1,4-benzodiazepin-2-one, **7a**
- 7,8-dimethoxy-1-methyl-5-[3-(4-phenyl-1, 3-thiazol-2-yl)phenyl]-1,3-dihydro-2H-1,4-benzodiazepin-2-one, **9b** <u>and</u>

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7,8-dimethoxy-l-methyl-5-(3-pyridyl)-1,3-dihydro-1,4-benzodiazepin-2-one, 17b.

49. (Currently Amended) A pharmaceutical composition comprising at least one compound such as defined In in accordance with claim 36 and a pharmaceutically acceptable vehicle or excipient.

Claims 50-54 cancelled.

55. (New) A compound represented by formula (1) according to claim 39, wherein R5 is a phenyl group substituted in positions 3 and 4 by a methylenedioxy chain that forms a ring with said phenyl group.

56. (New) A compound represented by formula (1)

$$R_{6}$$
 $R_{7}$ 
 $R_{6}$ 
 $R_{6}$ 
 $R_{7}$ 
 $R_{7}$ 
 $R_{7}$ 
 $R_{8}$ 
 $R_{1}$ 
 $R_{2}$ 
 $R_{3}$ 
 $R_{3}$ 
 $R_{3}$ 
 $R_{4}$ 
 $R_{5}$ 
 $R_{5}$ 

in which:

Z represents oxygen, sulfur or -NR<sub>2</sub>;

 $R_1$  is selected from the group consisting of hydrogen,  $(C_1-C_6)$  alkyl,  $(C_6-C_{18})$  aryl,-  $(C_1-C_6)$  alkyl $(C_6-C_{18})$  aryl and  $(C_6-C_{18})$  aryl $(C_1-C_4)$  alkyl;

 $R_2$  is selected from the group consisting of hydrogen, ( $C_1$ - $C_6$ ) alkyl, ( $C_6$ - $C_{18}$ )aryl, ( $C_1$ - $C_6$ )alkyl( $C_6$ - $C_{18}$ )aryl and ( $C_6$ - $C_{18}$ )aryl( $C_1$ - $C_4$ )aryl, and wherein Z is -NR2,  $R_1$  and  $R_2$  taken together may form a linear- or branched-hydrocarbon chain having from 2 to 6 carbon atoms, optionally containing one or more double bonds and optionally containing oxygen, sulfur or nitrogen;

 $R_3$  and  $R_3$ ', are independently selected from the group consisting of hydrogen, -( $C_1C_{12}$ ) alkyl, ( $C_3$ - $C_6$ ) cycloalkyl, ( $C_6$ - $C_{18}$ ) aryl, ( $C_6$ - $C_{18}$ )aryl( $C_1$ - $C_4$ )alkyl, ( $C_1$ - $C_{12}$ )alkyl( $C_6$ - $C_{18}$ )aryl, ( $C_5$ - $C_{18}$ ) heterocycle, aromatic or not, containing 1 to 3 heteroatoms, -NO<sub>2</sub>, CF<sub>3</sub>, CN, NR'R", SR', OR', COOR' and CONR'R", wherein R' and R" are independently selected from the group consisting of hydrogen, ( $C_1$ - $C_6$ ) alkyl, ( $C_3$ - $C_6$ ) cycloalkyl, ( $C_6$ - $C_{12}$ ) aryl and a ( $C_5$ - $C_{12}$ ) heterocycle, aromatic or not, containing 1 to 3 heteroatoms;

R<sub>5</sub> represents a phenyl group substituted at least in position 3; a naphthyl group; a (C<sub>5</sub>-C<sub>18</sub>) heterocycle, aromatic or not, containing 1 to 3 heteroatoms, selected from the group consisting of pyridyl, isoquinolyl, quinolyl and piperazinyl, provided that, when R<sub>5</sub> is a naphthyl group substituted in position 6, said naphthyl group is not attached to the rest of the molecule at position 2, or when R<sub>5</sub> is a pyridyl group, said pyridyl group is not attached to the rest of the molecule at position 4, or when R<sub>5</sub> is a tetrahydro-1,2,3,4isoquinolyl group, said group is not attached to the rest of the molecule at position 2, or when R<sub>5</sub> represents a phenyl group substituted at least in position 3, said substituent is selected from the group consisting of: alkyl, halogenoalkyl, cycloalkyl, alkenyl, alkynyl, aralkyl, aryl, (C<sub>5</sub>-C<sub>18</sub>) heterocycle, aromatic or not, hydroxy, = O, NO<sub>2</sub>, NH<sub>2</sub>, CN, COR', COOR', (C<sub>1</sub>-C<sub>6</sub>)alkoxy, (di)(C<sub>1</sub>-C<sub>6</sub>)alkylamino, NHCOR', CONR'R" group, in which R' and R" are as defined hereinabove, CHO, CONH2 and phenyl optionally substituted by an acetyl group, halogen, CONH2 or CN; prop-1-ynyl optionally substituted by a benzyloxy or tert-butyl carbamate; hex-1-ynyl optionally substituted by CN or NH2; pentyl optionally substituted by CONH2; hexyl; piperidinyl optionally substituted by a prop-1ynyl, benzylaminomethyl, acetamide, aminomethyl, NH2CS-, 4-phenyl-1,3-thiazol-2-yl, -CONHBenzyl, -COOEthyl, -CONHiPropyl, -CONH-(CH2)n-CONH2 (wherein n represents a whole number from 1 to 6), a -CONR'R" group, wherein R' and R" are independently C1-C6 alkyl or hydrogen, -(4-benzylpyperazin-1-yl)carbonyl, -CONH-(CH2)n-phenyl (wherein n represents a whole number from 1 to 6), imidazolyl, piperazinyl optionally substituted by a phenyl group, or R5 is substituted in positions 3 and 4 by halogens or by a hydrocarbon chain, optionally containing a heteroatom, thereby forming a ring with the phenyl;

 $R_7$  and  $R_8$  are independently selected from the group consisting of hydrogen, halogen and a  $OR_{10}$ , group in which  $R_{10}$  represents hydrogen,  $(C_1-C_6)$  alkyl,  $(C_3-C_6)$  cycloalkyl,  $(C_6-C_{12})$  aryl, or a  $(C_5-C_{12})$  heterocycle, aromatic or not, containing 1 to 3 heteroatoms, wherein at least one  $R_7$  and  $R_8$  represents a  $OR_{10}$  group as defined hereinabove;

 $R_6$  and  $R_9$  are independently selected from the group consisting of hydrogen, halogen, alkyl, cycloalkyl, alkenyl, alkynyl, aryl, aralkyl, heterocycle, aromatic or not, and an  $OR_{10}$  group, wherein R10 is as defined hereinabove;

further wherein the alkyl, cycloalkyl, alkenyl, alkynyl, aralkyl, aryl and heterocycle, aromatic or not, groups, or the hydrocarbon chain defined hereinabove are optionally substituted by one or more substituents independently selected from the group consisting of halogen, alkyl, halogenoalkyl, cycloalkyl, alkenyl, alkynyl, aralkyl, aryl, heterocycle, aromatic or not, hydroxy, =O, NO<sub>2</sub>, NH<sub>2</sub>, CN, COR', COOR', (C<sub>1</sub>-C<sub>6</sub>)alkoxy, (di)(C<sub>1</sub>-C<sub>6</sub>)alkylamino, NHCOR' and CONR'R" groups, in which R' and R" are as defined hereinabove,

and pharmaceutically acceptable salts thereof,

with the exception of compounds represented by formula (1) in which:

R1 is alkyl, R3, R3', R6 and R9 are hydrogen, and R5 is a phenyl group substituted at least in position 3 by a methoxy group;

R1 is alkyl or hydrogen, R3, R3', R6 and R9 are hydrogen, R5 is a phenyl group substituted only in position 3 by chlorine or bromine;

R1 is alkyl, R3, R3', R6 and R9 are hydrogen, R5 is a phenyl group substituted at least in position 3 by a CH2OH group;

R1, R3, R3', R6 and R9 are hydrogen, R5 is a phenyl group substituted only in position 3 by CF3,

R1 is alkyl, R3, R3', R6 and R9 are hydrogen, R5 is a phenyl group substituted in positions 3 and 5 by CF3,

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R1 is alkyl, R3, R3', R6 and R9 are hydrogen, R7 and R8 are methoxy, R5 is a phenyl group substituted in position 3 by a phenyl group, and

R1 is alkyl, R3, R3', R6 and R9 are hydrogen, R7 and R8 are methoxy, R5 is a phenyl group substituted in position 3 by a phenylethynyl group.